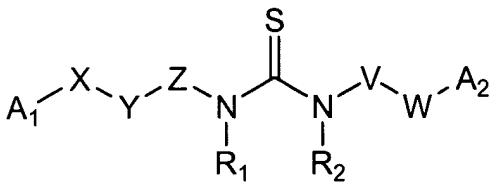
AMENDMENTS TO THE CLAIMS

1. (Original) A compound of Formula 1



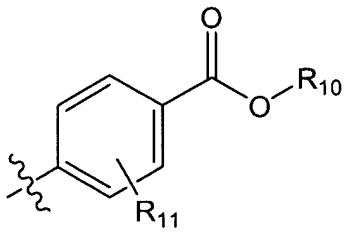
Formula 1

or a pharmaceutically acceptable salt thereof, wherein

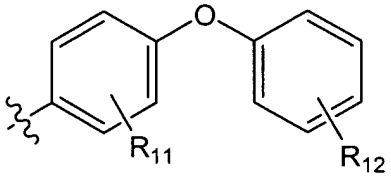
A_1 is an optionally substituted di-alkylamino, an optionally substituted aryl group, an optionally substituted 5- or 6- membered heteroaryl group, an optionally substituted bicyclic heteroaryl group having a 5-membered heteroaryl ring fused to a phenyl ring, an optionally substituted partially unsaturated or aromatic heterocyclic group having two 6-membered rings, an optionally substituted 5- to 7-membered heterocycloalkyl group containing at least one nitrogen atom and 0 or 1 additional heteroatoms, an optionally substituted partially unsaturated 5- to 7-membered heterocycloalkyl group containing at least one nitrogen atom and 0 or 1 additional heteroatoms, a 5- or 6-membered heterocycloalkyl group fused to a phenyl or heteroaryl ring, or a fused or spiro 8 to 11-membered bicyclic heterocycloalkyl group containing at least one nitrogen atom and 0 to 3 additional heteroatoms;

A_2 is

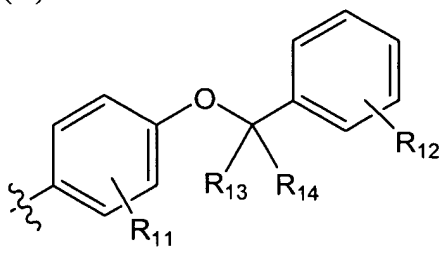
(i)



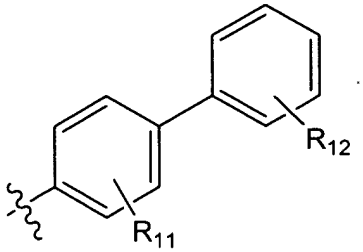
(ii)



(iii)



(iv)



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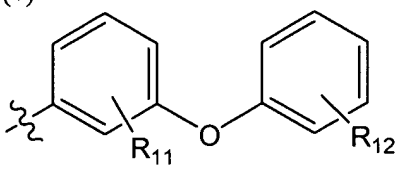
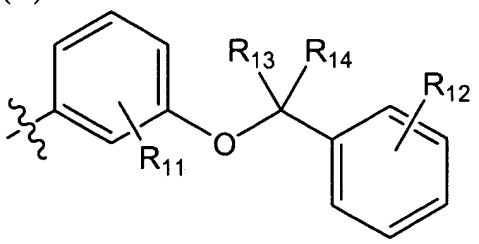


(vi)



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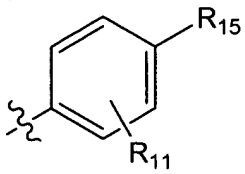
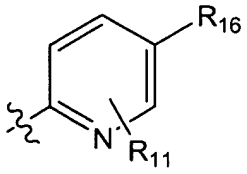


(viii)



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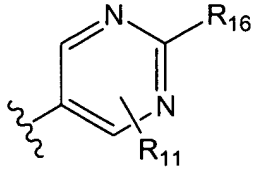
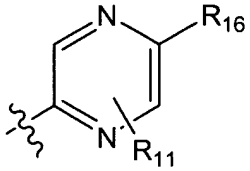
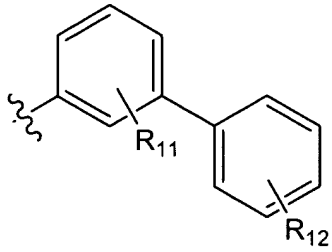


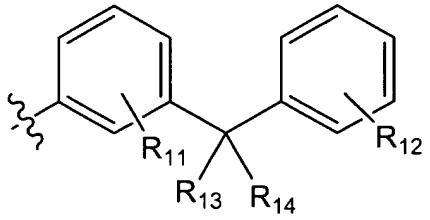
(x)



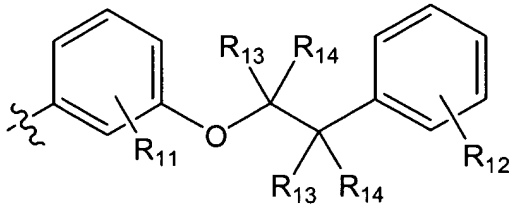
(xi)



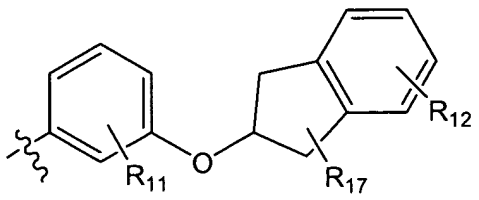
(xii)



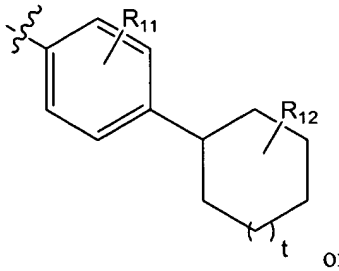
(xiii)



(xiv)

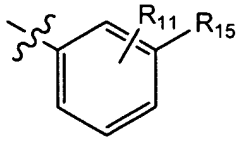


(xv)



or

(xvi)



;

t is 0 or 1;

X and W are independently O, S, NR, or absent, where R is hydrogen, optionally substituted C_1-C_6alkyl, or optionally substituted aryl(C_0-C_4alkyl);

V is C_1-C_6 alkyl, C_2-C_6alkenyl, C_3-C_7cycloalkyl, or absent;

Y is C_1-C_6 alkyl, C_1-C_6 alkyl substituted with C_3-C_7cycloalkyl, C_2-C_6alkenyl, C_3-C_7cycloalkyl, or absent;

wherein when V is absent, W is absent;

Z is carbonyl, thiocarbonyl, imino, or C_1-C_6alkylimino;

R_1 and R_2 are independently hydrogen, or

R_1 and R_2 are independently C_1-C_6alkyl, C_2-C_6 alkenyl, or C_2-C_6 alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkoxy, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, or

R_1 and R_2 are joined to form a 5- to 7-membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy;

R_{10} is C_1-C_6alkyl;

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1-C_6alkyl, C_1-C_6alkoxy, mono- and di-(C_1-C_6alkyl)amino, C_2-C_6alkanoyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl; and

R_{13} and R_{14} are independently chosen at each occurrence from hydrogen and C_1-C_4alkyl;

R_{15} is C_4-C_6alkoxy or C_4-C_6alkyl;

R_{16} is C_2-C_6alkoxy or C_2-C_6alkyl; and

R_{17} represents 0 to 2 substituents independently chosen from halogen, methyl, and methoxy;

2. (Original) A compound or salt according to Claim 1 wherein

A_1 is a di-(C_1-C_6alkyl)amino, an aryl group, a 5- or 6- membered heteroaryl group, a bicyclic heteroaryl group having a 5-membered heteroaryl ring fused to a phenyl ring, a partially unsaturated or aromatic heterocyclic group having two 6-membered rings, a 5- to 7-membered heterocycloalkyl group containing at least one nitrogen atom and 0 or 1 additional heteroatoms, a partially unsaturated 5- to 7-membered heterocycloalkyl group containing at least one nitrogen atom and 0 or 1 additional heteroatoms, a 5- or 6-membered heterocycloalkyl group fused to a phenyl or

heteroaryl ring, or a fused or spiro 8 to 11-membered bicyclic heterocycloalkyl group containing at least one nitrogen atom and 0 to 3 additional heteroatoms; each of which A_1 is substituted with 0 to 5 substituents independently chosen from:

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, $-CONH_2$, $-SO_2NH_2$, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy,

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl,
-mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl, and

(c) $-GR_a$ where

G is chosen from $-(CH_2)_n-$, C_2-C_4alkenyl, C_2-C_4alkynyl, $-(CH_2)_nO(CH_2)_m-$, and $-(CH_2)_nN(CH_2)_m-$, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl;

X and W are independently O, S, NR, or absent,

where R is hydrogen or R is C_1-C_6alkyl or aryl(C_0-C_4alkyl), each of which is substituted with 0 to 5 substitutents independently chosen from halogen, hydroxy, cyano, amino, nitro, oxo, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, C_1-C_6alkyl, C_1-C_6alkoxy, and mono- and di-(C_1-C_6alkyl)amino;

V is independently C_1-C_6 alkyl, C_2-C_6alkenyl, C_3-C_7cycloalkyl, or absent;

Y is C_1-C_6 alkyl, C_1-C_6 alkyl substituted with C_3-C_7cycloalkyl, C_2-C_6alkenyl, C_3-C_7cycloalkyl, or absent;

wherein when V is absent, W is absent;

Z is carbonyl, thiocarbonyl, or imino; and

R_1 and R_2 are independently hydrogen, or

R_1 and R_2 are independently C_1-C_6alkyl, C_2-C_6 alkenyl, or C_2-C_6 alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkoxy, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, or

R_1 and R_2 are joined to form a 5- to 7-membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-

unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

Claims 3-5 (Canceled)

6. (Currently Amended) A compound or salt according to Claim ~~1 or~~ 2 in which Z is carbonyl.

7. (Currently Amended) A compound or salt according to ~~any one of~~ Claims ~~1 to~~ 6 in which X is oxygen and Y is $-CH_2-$.

Claim 8 (Canceled)

9. (Currently Amended) A compound or salt according to ~~any one of~~ Claims ~~1 to~~ 6 wherein when X and Y are absent.

10. (Currently Amended) A compound or salt according to ~~any one of Claims 1 to 9~~ Claim 6 wherein V and W are absent.

Claims 11-12 (Canceled).

13. (Currently Amended) A compound or salt according to Claim ~~12~~ 6 in which R_1 and R_2 are independently hydrogen, methyl, or ethyl.

14. (Original) A compound or salt according to Claim 13 in which R_1 and R_2 are both hydrogen.

Claims 15-16 (Canceled).

17. (Currently Amended) A compound or salt according to ~~any one of Claims 2 to 16~~ Claim 6 wherein

A_1 is aryl, a partially unsaturated heterocyclic group, or heteroaryl group;

substituted with 0 to 5 substituents independently chosen from:

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -$CONH_2$, -SO_2NH_2, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, and

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl, mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl, and

(c) –GR_a where

G is chosen from -$(CH_2)_n$-, C_2-C_4alkenyl, C_2-C_4alkynyl, -$(CH_2)_nO(CH_2)_m$-, and -$(CH_2)_nN(CH_2)_m$-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

18. (Original) A compound or salt according to Claim 17 wherein

A_1 is phenyl, naphthyl, pyridyl, pyrimidinyl, thienyl, pyrrolyl, furanyl, pyrazolyl, imidazolyl, thiazolyl, triazolyl, thiadiazolyl, oxazolyl, isoxazolyl, benzofuranyl, benzothiazolyl, benzothiophenyl, benzoxadiazolyl, benzo[*d*]oxazolyl, dihydrobenzodioxynyl, indolyl, pyrazolopyrimidinyl, thienylpyrazolyl, benzopyranyl, or 4*H*-chromenyl,

each of which is substituted with 0 to 5 substituents independently chosen from

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -$CONH_2$, -SO_2NH_2, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, and

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl, -mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl, and

(c) –GR_a where

G is chosen from -$(CH_2)_n$-, C_2-C_4alkenyl, C_2-C_4alkynyl, -$(CH_2)_nO(CH_2)_m$-, and

$-(CH_2)_nN(CH_2)_m-$, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

Claim 19 (Canceled)

20. (Currently Amended) A compound or salt according to Claim ~~19~~ 18 in which A_1 is substituted with 0 to 5 substituents independently chosen from

(a) halogen, hydroxy, cyano, amino, nitro, oxo, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, and

(b) C_1-C_6alkyl, C_1-C_6alkoxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_4)alkyl, mono- and di-(C_1-C_4alkyl)amino, and mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl; and

(c) $-GR_a$ where

G is from $-(CH_2)_n-$, $-(CH_2)_nO(CH_2)_m-$, and $-(CH_2)_nN(CH_2)_m-$, and

R_a is C_3-C_8cycloalkyl, 5 or 6-membered heterocycloalkyl containing 1 or 2 heteroatoms independently chosen from O, S, and N, 5- or 6-membered heteroaryl containing 1, 2, or 3 heteroatoms independently chosen from O, S, and N, indanyl, and phenyl,

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_2alkyl)amino, and C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

21. (Original) A compound or salt according to Claim 20 in which

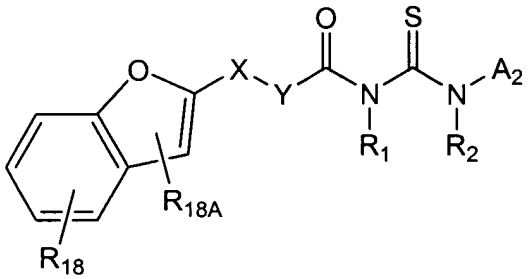
A_1 is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, oxo, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, C_1-C_6alkyl, C_1-C_6alkoxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_4)alkyl, mono- and di-(C_1-C_4alkyl)amino, and mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl.

Claims 22-39 (Canceled).

40. (Currently Amended) A compound or salt according to ~~any one of Claims 1, 2, or 24 to 38,~~ Claim 2 wherein

A_1 is pyrazinyl, pyridyl, or quinaxolinyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

41. (Currently Amended) A compound or salt according to ~~any one of Claims 1, 2, or 24 to 38~~ Claim 2 of Formula 17, wherein



Formula 17

wherein

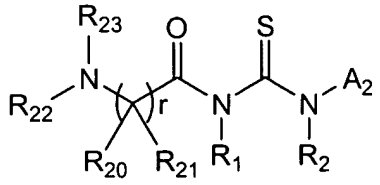
R_{18A} is hydrogen, halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, or C_1-C_2haloalkoxy; and

R_{18} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

42. (Original) A compound or salt according to Claim 41 in which

X and Y are absent; and R_1 and R_2 are independently hydrogen or methyl.

43. (Currently Amended) A compound or salt according to ~~any one of Claims 1, 2, or 24 to 38~~ Claim 2 of Formula 18, wherein



Formula 18

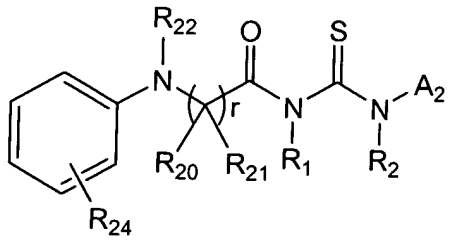
r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1-C_4alkyl; or R_{20} and R_{21} are joined to form a C_3-C_7cycloalkyl group; and

R_{22} and R_{23} are independently chosen C_1-C_6 alkyl groups; each of which is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy,

mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

44. (Currently Amended) A compound or salt according to ~~any one of Claims 1, 2, or 24 to 38~~ Claim 2 of Formula 19, wherein



Formula 19

r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1-C_4alkyl; or R_{20} and R_{21} are joined to form a C_3-C_7cycloalkyl group;

R_{22} is C_1-C_6 alkyl which is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

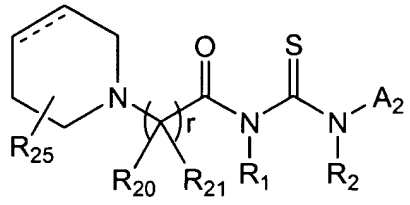
R_{24} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, -COOH, -$CONH_2$, -SO_2NH_2, -SH, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl,

-mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl.

45. (Original) A compound or salt according to Claim 44, wherein

R_{24} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

46. (Currently Amended) A compound or salt according ~~any one of Claims 1, 2 or 24 to 38~~ to Claim 2 of Formula 20, wherein



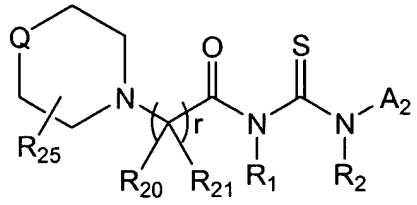
Formula 20

r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1-C_4alkyl; or R_{20} and R_{21} are joined to form a C_3-C_7cycloalkyl group; and

R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

47. (Currently Amended) A compound or salt according ~~any one of Claims 1, 2 or 24 to 38~~ to Claim 2 of Formula 21, wherein



Formula 21

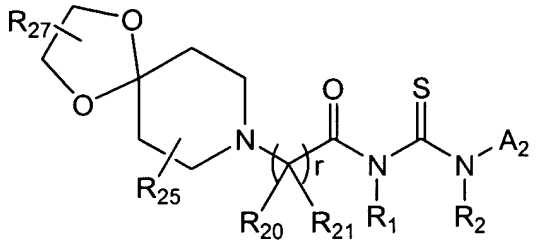
r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1-C_4alkyl; or R_{20} and R_{21} are joined to form a C_3-C_7cycloalkyl group; and

R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

Q is O, S, or NR_{26}; where R_{26} is hydrogen or R_{26} is C_1-C_6alkyl, phenyl, pyridyl, or pyrimidinyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_6alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

48. (Currently Amended) A compound or salt according ~~any one of Claims 1, 2 or 24 to 38~~ to Claim 2 of Formula 22, wherein



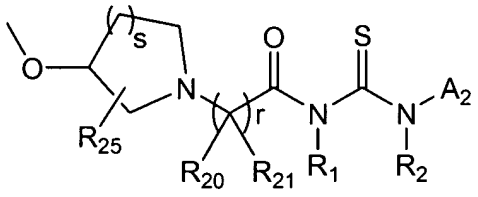
Formula 22

r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1-C_4alkyl; or R_{20} and R_{21} are joined to form a C_3-C_7cycloalkyl group; and

R_{25} and R_{27} each represent 0 to 2 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

49. (Currently Amended) A compound or salt according ~~any one of Claims 1, 2 or 24 to 38~~ to Claim 2 of Formula 23, wherein



Formula 23

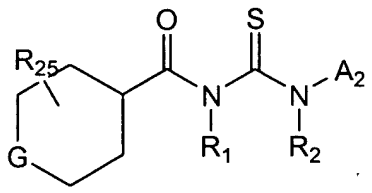
r is 1, 2, or 3;

s is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1-C_4alkyl; or R_{20} and R_{21} are joined to form a C_3-C_7cycloalkyl group; and

R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

50. (Currently Amended) A compound or salt according ~~any one of Claims 1, 2 or 24 to 38~~ to Claim 2 of Formula 24, wherein



Formula 24

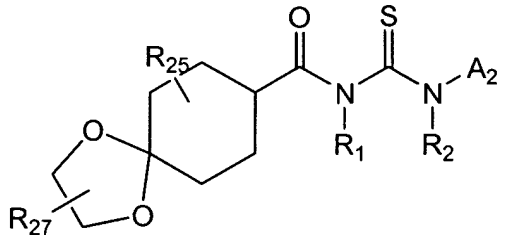
R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_6alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

G is O, S, SO_2, or NR_{26}; where R_{26} is hydrogen or R_{26} is C_1-C_6alkyl, phenyl, pyridyl, or pyrimidinyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_6alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

51. (Original) A compound or salt according to Claim 50, wherein

R_{25} represents a di-(C_1-C_6alkyl)amino substituent and 0 to 2 additional substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_6alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

52. (Currently Amended) A compound or salt according to ~~any one of Claims 1, 2, or 24 to 38~~ Claim 2 of Formula 25, wherein



Formula 25

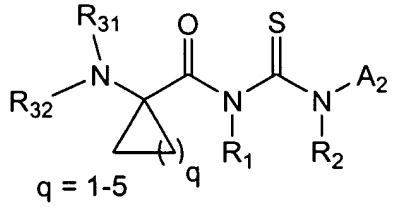
R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{27} represents 0 to 2 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

53. (Original) A compound or salt according to Claim 52, wherein

R_{25} represents a di-(C_1-C_6alkyl)amino substituent and 0 to 2 additional substituents independently chosen from halogen, hydroxy, cyano, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_6alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

54. (Currently Amended) A compound or salt according ~~any one of Claims 1, 2 or 24 to 38~~ to Claim 2 of Formula 26, wherein

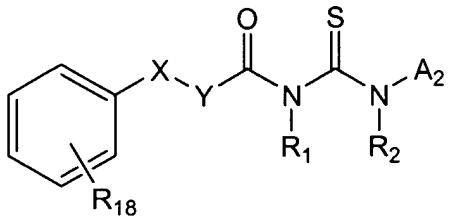


Formula 26

q is an integer from 1 to 5; and

R_{31} and R_{32} are independently chosen from C_1-C_6alkyl and phenyl; each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

55. (Currently Amended) A compound or salt according to ~~any one of Claims 1, 2, or 24 to 38~~ Claim 2 of Formula 27, wherein



Formula 27

wherein

R_{18} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

56. (Original) A compound or salt according to Claim 55 wherein X is oxygen and Y is $-CH_2-$.

57. (Original) A compound or salt according to Claim 55 wherein X and Y are absent.

58. (Currently Amended) A compound or salt according to ~~any one of Claims 1, 2, or 24 to 38~~ Claim 2, wherein

A_1 is 5-membered heteroaryl group selected from furan-2-yl, furan-3-yl, isoxazol-3-yl, isoxazol-4-yl, thiophen-2-yl, thiophen-3-yl, pyrrol-2-yl, pyrrol-3-yl, and pyrazolyl; each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_2-C_4alkenyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

59. (Original) A compound or salt according to Claim 58 wherein X is oxygen and Y is $-CH_2-$.

60. (Original) A compound or salt according to Claim 58 wherein X and Y are absent.

61. (Currently Amended) A compound or salt according to ~~any one of Claims 1, 2, or 24 to 38~~

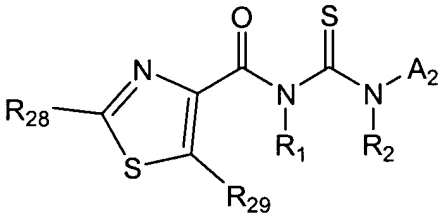
Claim 2, wherein

A_1 is pyridin-2-yl or pyridin-3-yl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

62. (Original) A compound or salt according to Claim 61 wherein X is oxygen and Y is $-CH_2-$.

63. (Original) A compound or salt according to Claim 61 wherein X and Y are absent.

64. (Currently Amended) A compound or salt according to ~~any one of Claims 1, 2, or 24 to 38~~ Claim 2 of Formula 28, wherein



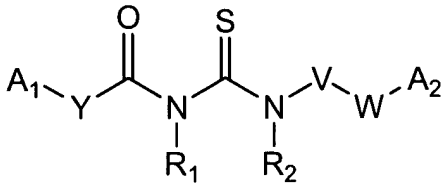
Formula 28

wherein

R_{28} is phenyl or pyridyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; and

R_{29} is hydrogen, methyl or ethyl.

65. (Original) A compound or pharmaceutically acceptable salt of Formula IB, wherein



Formula IB

A_1 is di-(C_1-C_8alkyl)amino, an *N*-(C_1-C_6alkyl)-*N*-phenyl-amino group, an *N*-(C_1-C_6alkyl)-*N*-pyridyl amino group, a 5- to 7-membered monocyclic heterocycloalkyl group covalently bound to a point of attachment in Formula IB via a Nitrogen atom, a 5- to 7-membered monocyclic partially unsaturated heterocyclic group covalently bound to a point of attachment in Formula IB via a Nitrogen atom, a 5- to 7- membered heterocycloalkyl group covalently bound to a point of attachment in Formula IB via a Carbon atom which is adjacent to a Nitrogen atom, or an 8- to 11-membered bicyclic heterocycloalkyl in which the rings are fused or spiro covalently bound to a

point of attachment in Formula IB via a Nitrogen atom;

A_2 is C_3-C_8 cycloalkyl, a partially unsaturated or aromatic carbocyclic group, or a saturated, partially unsaturated, or aromatic heterocyclic group;

each of which A_1 and A_2 is substituted with 0 to 5 substituents independently chosen from (a), (b), and (c), where

(a) is independently chosen from halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -$CONH_2$, -SO_2NH_2, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy,

(b) is independently chosen from C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl, -mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl, and

(c) is –GR_a where G is chosen from -$(CH_2)_n$-, C_2-C_4alkenyl, C_2-C_4alkynyl, -$(CH_2)_nO(CH_2)_m$-, and -$(CH_2)_nN(CH_2)_m$-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl;

W is O, S, NR, or absent, where R is hydrogen or R is C_1-C_6alkyl or aryl(C_0-C_4alkyl), each of which is substituted with 0 to 5 substitutents independently chosen from halogen, hydroxy, cyano, amino, nitro, oxo, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, C_1-C_6alkyl, C_1-C_6alkoxy, and mono- and di-(C_1-C_6alkyl)amino;

V is C_1-C_6 alkyl, C_2-C_6alkenyl, C_3-C_7cycloalkyl, or absent; and when V is absent, W is absent;

Y is C_1-C_6 alkyl substituted with

0 or 1 of C_3-C_7cycloalkyl, a 5- to 7-membered monocyclic heterocycloalkyl, or 8- to 11- membered bicyclic heterocycloalkyl in which the rings are fused or spiro; each of which substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy; or

Y is absent;

R_1 and R_2 are independently hydrogen or C_1-C_6alkyl, C_2-C_6 alkenyl, or C_2-C_6 alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkoxy, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, or R_1 and R_2 are joined to form a 5- to 7-

membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy.

66. (Original) A compound or salt according to Claim 65 wherein V and W are absent.

67. (Original) A compound or salt according to Claim 65 in which Y is absent.

Claims 68-70 (Canceled)

71. (Currently Amended) A compound or salt according to Claim ~~70~~ 67 in which R_1 and R_2 are independently hydrogen or methyl.

72. (Currently Amended) A compound or salt according to ~~any one of Claims 65 to 71~~ Claim 71 wherein

A_2 is C_5-C_7cycloalkyl, phenyl, pyridyl, naphthyl, pyrimidinyl, pyrazinyl, benzothiazolyl, benzodioxyl, quinolinyl, or isoquinolinyl, each of which is substituted with 0 to 5 substituents independently chosen from (a), (b), and (c) where

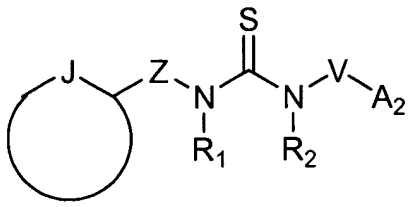
(a) is chosen from halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -$CONH_2$, -SO_2NH_2, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy,

(b) is chosen from C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkanoyl, and C_1-C_8alkoxycarbonyl, and

(c) is $-GR_a$ where G is chosen from -$(CH_2)_n$-, C_2-C_4alkenyl, C_2-C_4alkynyl, -$(CH_2)_nO(CH_2)_m$-, and -$(CH_2)_nN(CH_2)_m$-, where n and m are independently 0, 1, 2, 3, or 4; and R_a is chosen from C_3-C_8cycloalkyl, piperidinyl, piperazinyl, morpholinyl, tetrahydroisoquinolinyl, indanyl, tetrahydronaphthyl, phenyl, pyridyl, benzothiophenyl, and benzofuranyl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

Claim 73. (Original) A compound of Formula II



Formula II

or a pharmaceutically acceptable salt thereof wherein

A_2 is C_3-C_8 cycloalkyl, a partially unsaturated or aromatic carbocyclic group, a saturated, partially unsaturated, or an aromatic heterocyclic group substituted with 0 to 5 substituents independently chosen from:

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -$CONH_2$, -SO_2NH_2, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, and

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkenyloxy, C_1-C_4alkoxy(C_1-C_4alkyl), amino(C_1-C_6)alkyl, mono- and di-(C_1-C_6alkyl)amino, mono- and di-(C_1-C_4alkyl)aminoC_1-C_4alkyl, C_2-C_6alkanoyl, C_2-C_8alkanoyloxy, C_1-C_8alkoxycarbonyl,
-mono- and di-(C_1-C_6alkyl)carboxamide, (C_3-C_7cycloalkyl)carboxamide, mono- and di-(C_1-C_6alkyl)sulfonamide, C_1-C_6alkylthio, aryl(C_0-C_4alkyl)thio, C_1-C_6alkylsulfinyl, and C_1-C_6alkylsulfonyl, and

(c) –GR_a where

G is chosen from -$(CH_2)_n$-, C_2-C_4alkenyl, C_2-C_4alkynyl, -$(CH_2)_nO(CH_2)_m$-, and -$(CH_2)_nN(CH_2)_m$-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl;

V is C_1-C_6 alkyl, C_2-C_6alkenyl, or absent; and

Z is carbonyl, thiocarbonyl, or imino;

R_1 and R_2 are independently

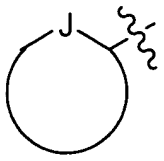
hydrogen, or

C_1-C_6alkyl, C_2-C_6 alkenyl, or C_2-C_6 alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkoxy, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, or

R_1 and R_2 are joined to form a 5- to 7-membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-

unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy;

the group:



is a group of Formula (i)

that is a saturated, partially unsaturated, or aromatic heterocyclic group where J is O, S, or NR_3 substituted with 0 to 5 substituents independently chosen from: (a), (b), and (c) above; and

R_3 is

(d) hydrogen, C_1-C_2haloalkyl, or C_1-C_2haloalkoxy;

(e) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_4alkoxy(C_1-C_4alkyl), or amino(C_1-C_6)alkyl, or

(f) $-LR_b$ where

L is chosen from $-(CH_2)_r-$, C_2-C_4alkenyl, C_2-C_4alkynyl, $-(CH_2)_rO(CH_2)_s-$, and $-(CH_2)_rN(CH_2)_s-$, where r and s are independently 0, 1, 2, 3, or 4; and

R_b is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (e) and (f) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

74. (Original) A compound or salt according to Claim 73 wherein Z is carbonyl.

75. (Currently Amended) A compound or salt according to ~~Claim 73 or~~ Claim 74 wherein V is absent or V is C_1-C_4alkyl.

76. (Currently Amended) A compound or salt according to ~~any one of Claims 73 to 75~~ Claim 75 wherein R_1 and R_2 are independently hydrogen or methyl.

77. (Currently Amended) A compound or salt according to ~~any one of Claims 73 to 76~~ Claim 76 wherein

A_2 is C_5-C_7cycloalkyl, phenyl, pyridyl, naphthyl, pyrimidinyl, pyrazinyl, benzothiazolyl, benzodioxyl, quinolinyl, or isoquinolinyl, each of which is substituted with 0 to 5 substituents independently

chosen from (a), (b), and (c).

78. (Currently Amended) A compound or salt according to ~~any one of Claims 73 to 77~~ Claim 77 wherein

A_2 is C_5-C_7cycloalkyl, phenyl, pyridyl, naphthyl, benzothiazolyl, benzodioxyl, quinolinyl, or isoquinolinyl, each of which is substituted with 0 to 5 substituents independently chosen from

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, $-CONH_2$, $-SO_2NH_2$, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy,

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkanoyl, and C_1-C_8alkoxycarbonyl, and

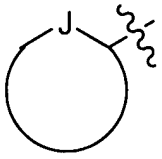
(c) $-GR_a$ where

G is chosen from $-(CH_2)_n-$, C_2-C_4alkenyl, C_2-C_4alkynyl, $-(CH_2)_nO(CH_2)_m-$, and $-(CH_2)_nN(CH_2)_m-$, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, piperidinyl, piperazinyl, morpholinyl, tetrahydroisoquinolinyl, indanyl, tetrahydronaphthyl, phenyl, pyridyl, benzothiophenyl, and benzofuranyl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

79. (Currently Amended) A compound or salt according to ~~any one of Claims 73 to 78~~ Claim 78 wherein



is a group of Formula (i)

where Formula (i) is a heteroaryl group that is pyridyl, pyrimidinyl, thienyl, pyrrolyl, furanyl, pyrazolyl, imidazolyl, thiazolyl, triazolyl, thiadiazolyl, oxazolyl, isoxazolyl, benzofuranyl, benzothiazolyl, benzothiophenyl, benzoxadiazolyl, benzo[*d*]oxazolyl, dihydrobenzodioxynyl, indolyl, pyrazolopyrimidinyl, or thienylpyrazolyl oriented such that the heteroatom J is adjacent to the point of attachment of the group of Formula (i)

the group of Formula (i) is substituted with 0 to 5 substituents independently chosen from: (a), (b), and (c);

J is S, O, or NR_3; and

R_3 is

(d) hydrogen, C_1-C_2haloalkyl, or C_1-C_2haloalkoxy;

(e) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_4alkoxy(C_1-C_4alkyl), or amino(C_1-C_6)alkyl, or

(f) $-LR_b$ where

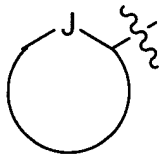
L is chosen from $-(CH_2)_r-$, C_2-C_4alkenyl, C_2-C_4alkynyl, $-(CH_2)_rO(CH_2)_s-$, and $-(CH_2)_rN(CH_2)_s-$, where r and s are independently 0, 1, 2, 3, or 4; and

R_b is chosen from C_3-C_8cycloalkyl, C_2-C_7monocyclic heterocycloalkyl, C_5-C_{10}bicyclicheterocycloalkyl, indanyl, tetrahydronapthyl, aryl, and heteroaryl;

each of which (e) and (f) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

80. (Original) A compound or salt of Claim 79 wherein

wherein



is a group of Formula (i)

where Formula (i) is a heteroaryl group that is pyridyl, pyrimidinyl, thienyl, pyrrolyl, furanyl, pyrazolyl, imidazolyl, thiazolyl, triazolyl, thiadiazolyl, oxazolyl, isoxazolyl, benzofuranyl, benzothiazolyl, benzothiophenyl, benzoxadiazolyl, benzo[*d*]oxazolyl, dihydrobenzodioxynyl, indolyl, pyrazolopyrimidinyl, or thienylpyrazolyl oriented such that the heteroatom J is adjacent to the point of attachment of the group of Formula (i) in Formula II;

the group of Formula (i) is substituted with 0 to 5 substituents independently chosen from:

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, $-CONH_2$, $-SO_2NH_2$, -SH, C_1-C_2haloalkyl, and C_1-C_2haloalkoxy, and

(b) C_1-C_6alkyl, C_2-C_6alkenyl, C_2-C_6alkynyl, C_1-C_6alkoxy, C_2-C_6alkanoyl, and C_1-C_8alkoxycarbonyl,

(c) $-GR_a$ where

G is chosen from $-(CH_2)_n-$, C_2-C_4alkenyl, C_2-C_4alkynyl, $-(CH_2)_nO(CH_2)_m-$, and $-(CH_2)_nN(CH_2)_m-$, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3-C_8cycloalkyl, piperidinyl, piperazinyl, morpholinyl, tetrahydroisoquinolinyl, indanyl, tetrahydronaphthyl, phenyl, pyridyl, benzothiophenyl, and benzofuranyl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_2-C_4alkanoyl, C_1-C_4alkoxycarbonyl, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl;

J is S, O, or NR_3, and

R_3 is

(d) hydrogen,

(e) C_1-C_6alkyl, or

(f) $-LR_b$ where

L is chosen from $-(CH_2)_r-$, $-(CH_2)_rO(CH_2)_s-$, and $-(CH_2)_rN(CH_2)_s-$, where r and s are independently 0, 1, 2, 3, or 4; and

R_b is chosen from C_3-C_8cycloalkyl, piperidinyl, piperazinyl, morpholinyl, indanyl, tetrahydronapthyl, phenyl, and pyridyl;

each of which (e) and (f) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1-C_4alkyl, C_1-C_4alkoxy, mono- and di-(C_1-C_4alkyl)amino, C_1-C_2haloalkyl, C_1-C_2haloalkoxy, and phenyl.

81. (Original) A compound or pharmaceutically acceptable salt thereof, in which the compound is selected from

1-(Furan-2-carbonyl)-3-(4-benzo[*d*]thiazol-2-yl-phenyl)-thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-[5-(benzo[*d*]oxazol-2-yl)-2-methyl]phenylthiourea;

1-(3-(Benzo[*d*]thiazol-2-yl)phenyl)-3-(2-phenoxyacetyl) thiourea;

1-(4-(Benzo[*d*]oxazol-2-yl)phenyl)-3-propionylthiourea;

1-(Pyridin-3-carbonyl)-3-(4-benzo[*d*]thiazol-2-yl-phenyl)-thiourea;

1-[3-(2-chlorophenyl-5-methyl-isoxazol-4-yl)-carbonyl]-3-(4-isopropylphenyl)thiourea;

Butyl4-(3-(2-phenoxyacetyl) thioureido)benzoate;

Butyl 4-(3-acetylthioureido)benzoate;

Butyl 4-(3-(2-(3-chlorophenoxy) acetyl) thioureido)benzoate;

Butyl 4-(3-(3-phenoxypropanoyl) thioureido)benzoate;

Butyl 4-(3-(2-(naphthalen-3-yloxy)acetyl)thioureido)benzoate;

Butyl 4-(3-(benzofuran-2-yl-carbonyl)thioureido)benzoate;

Ethyl 2-(4-(3-(2-phenoxyacetyl)thioureido)phenyl)acetate;

Ethyl 4-(3-(2-phenoxyacetyl)thioureido)benzoate;

Butyl 4-(3-(2-methoxyacetyl) thioureido)benzoate;

Butyl 4-(3-(2-(2,4-dichlorophenoxy) acetyl)thioureido)benzoate;

Butyl 4-(3-(2-(4-tert-butylphenoxy) acetyl)thioureido)benzoate;

Butyl 4-(3-(2-(4-(benzyloxy) phenoxy)acetyl)thioureido)benzoate;

Butyl 4-(3-(2-(2-methoxyphenoxy) acetyl)thioureido)benzoate;

Butyl 4-(3-(2-(o-tolyloxy)acetyl)thioureido)benzoate;

Butyl 4-(3-(2-(2,4,6-trichlorophenoxy)acetyl)thioureido)benzoate;

Butyl 4-(3-(3,4-dichlorophenyl) carbonyl)thioureido)benzoate;

1-(3,4-dichlorophenyl-carbonyl)-3-(3-trifluromethylphenyl)thiourea;

1-(3,4-Dichlorophenyl-carbonyl)-3-(3-benzoxy-phenyl)thiourea;

1-(3,4-Dichlorophenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(3,4-Diflurophenyl-carbonyl)-3-(3-(5-methylfuran-2-yl)-phenyl)thiourea;

Butyl 4-(3-(naphth-2-yl) carbonyl)thioureido)benzoate;

Butyl 4-(3-(4-cyanophenyl) carbonyl)thioureido)benzoate;

Butyl 4-(3-(methylacetate) carbonyl)thioureido)benzoate;

1-((Benzofuran-2-yl-carbonyl)-3-(4-(phenoxy)-phenyl)thiourea;

Butyl 4-(3-(2-(3,4-dichlorophenoxy)acetyl)thioureido)benzoate;

1-(4-Butylphenyl)-3-(2-phenoxyacetyl)thiourea;

(Amino-(3-(benzyloxy)phenyl) methanethiocarbamoyl)methyl acetate;

1-(3-(Methylthio)propanoyl)-3-(3-(benzyloxy)phenyl)thiourea;

1-(2-(2-Chlorophenoxy)acetyl)-3-(3-(benzyloxy)phenyl)thiourea;

Butyl 4-(3-(naphth-1-yl) carbonyl)thioureido)benzoate;

(S)-1-(Amino-*N*-p-(butylacetate) methanethiocarbamoyl)ethyl acetate;

Butyl 4-(3-(2-(2-methoxyethoxy)acetyl)thioureido)benzoate;

(Amino*N*-(4-cyclohexylphenyl)methanethiocarbamoyl)(phenyl)methyl acetate;

Ethyl 3-(amino-*N*-(4-cyclohexylphenyl)methanethiocarbamoyl)propanoate;

1-Butyryl-3-(4-cyclohexylphenyl)thiourea;

(S)-1-(Amino-*N*-(4-cyclohexylphenyl)methanethiocarbamoyl)ethyl acetate;

1-(3-(Benzyloxy)phenyl)-3-(2-hydroxyacetyl)thiourea;

Butyl 4-(3-(2-(2,6-dichlorophenoxy) acetyl)thioureido)benzoate;

Butyl 4-(3-(2-(3-methoxyphenoxy) acetyl) thioureido)benzoate;

1-[(1-methyimidazol-2-yl) –carbonyl]-3-(3-benzoxy-phenyl)thiourea;

tert-Butyl 2-(aminoN-(3-(benzyloxy) phenyl)methanethiocarbamoyl) pyrrolidone-1-carboxylate;

Butyl 4-(3-(pyrrolidin-1-yl) carbonyl)thioureido)benzoate;

Butyl 4-(3-(1-methyl-benzofuran-2-yl) carbonyl)thioureido)benzoate;

1-(4-Hexylphenyl)-3-(2-phenoxyacetyl)thiourea;

1-(4-(Pentyloxy)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(4-pentyloxy)-phenyl)thiourea;

1-(4-Pentylphenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(4-pentyl)-phenyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(4-pentyloxy)-phenyl)thiourea;

1-(4-Butoxyphenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(3-phenyl)-phenyl)thiourea;

1-(2-Phenoxyacetyl)-3-(3-phenyl)-phenylthiourea;

Isopropyl 4-(3-(benzofuran-2-yl) carbonyl)thioureido)benzoate;

1-(2-phenoxyacetyl)-3-94-fluoro-phenyl)-phenylthiourea;

1-(3-benzylphenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(3-benzyl)-phenyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(4-benzyl)-phenyl)thiourea;

1-(4-(p-Tolyloxy)phenyl)-3-(2-phenoxyacetyl)thiourea;

Isobutyl 4-(3-(benzofuran-2-yl) carbonyl)thioureido)benzoate;

Isobutyl 4-(3-(2-phenoxyacetyl)thioureido)benzoate;

1-(2-(phenylmethanone)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-(3-(Phenylcarbamoyl)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-(3-(2-Methylpyrimidin-4-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-(4-(4-Chlorophenoxy)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(4-Chlorophenoxy)-phenyl)thiourea;

1-(4-(3,4-Dihydroisoquinolin-2(1H)-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-(3-Fluoro-4-(octahydroquinolin-1(2H)-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-(3-Fluoro-4-(octahydroquinolin-1(2H)-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-(3-Fluoro-4-(piperidin-1-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(3-fluoro-4-(piperidin-1-yl)phenyl)-phenyl)thiourea;

1-(3-(3-Methoxybenzyloxy) phenyl)-3-(2-phenoxyacetyl) thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(3-(3-methoxybenzyloxy)) phenyl)-phenyl)thiourea;

1-(3-(2-Methoxybenzyloxy) phenyl)-3-(2-phenoxyacetyl) thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(3-(2-Methoxybenzyloxy) phenyl)-thiourea;

1-(3-(4-Methoxybenzyloxy) phenyl)-3-(2-phenoxyacetyl) thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(3-(cyclohexylmethoxy) phenyl)-thiourea;

1-(3-(Cyclohexylmethoxy) phenyl)-3-(2-phenoxyacetyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(4-(5,6-dihydropyridin-1(2H)-yl))phenyl)-thiourea;

1-((5-Methoxy-benzofuran-2-yl-carbonyl)-3-(3-benzyloxy-phenyl)-thiourea;

Butyl 4-(3-(5-chloro-benzofuran-2-yl) carbonyl)thioureido)benzoate;

1-(7-Methoxy-benzofuran-2-yl-carbonyl)-3-(3-methoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(2,3,4-tetrahydronapthalen-2-yl-carbonyl)-3-(3-(3-methoxybenzyloxy))phenyl)-phenyl)thiourea;

1-(2-(4-(Trifluoromethoxy) phenoxy) acetyl)-3-(3-(benzyloxy)phenyl) thiourea;

1-(3-(Benzyloxy)phenyl)-3-(2-(pyridin-3-yloxy)acetyl) thiourea;

1-(4-oxo-4-*H*-chromen-2-yl-carbonyl)-3-(3-methoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(3-(Benzyloxy)phenyl)-3-(2-(pyridin-2-yloxy)acetyl) thiourea;

1-(Pyridin-2-yl-carbonyl)-3-(3-methoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(3-Chloro-benzo[*b*]thiophen-2-yl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(4-Trifluoromethoxy-phenyl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(5-Methylisoxazol-3-yl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(2-Methyl-5-phenyl-furan-3-yl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(4-Trifluoromethyl-phenyl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;

1-(3-Chloro-benzo[b]thiophen-2-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(4-Trifluoromethoxyphenyl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(3,5-Dimethylisoxazol-4-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(5-Methylisoxazol-3-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(4-Trifluoromethylphenyl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

(S)-1-(Amino-*N*-(3-(benzyloxy) phenyl)methanethiocarbamoyl)ethyl acetate;

(S)-1-(Amino-*N*-(3-phenoxyphenyl)methane thiocarbamoyl)ethyl acetate;

(S)-1-(Amino-*N*-(3-benzyl-phenyl)methanethiocarbamoyl)ethyl acetate;

Ethyl 1-(2-fluoro-4-(3-(benzofuran-2-yl-carbonyl)thioureido)phenyl)-4-phenylpiperidine-4-carboxylate;

1-(4-Trifluoromethylphenyl-carbonyl)-3-(2-phenylbenzo[d][1,3]dioxol-6-yl) phenyl)thiourea;

1-(3-Chloro-methylbenzo[b]thiophen-2-yl-carbonyl)-3-(2-phenylbenzo[d][1,3]dioxol-6-yl) phenyl)thiourea;

1-(4-Trifluoromethoxyphenyl-carbonyl)-3(2-phenylbenzo[d][1,3]dioxol-6-yl) thiourea;

1-(5-Methylisoxazol-3-yl-carbonyl)-3-(2-phenylbenzo[d][1,3]dioxol-6-yl) phenyl)thiourea;

1-(3-((R)-1-Phenylethoxy)phenyl)-3-(4-trifluoromethylphenyl-carbonyl)thiourea;

1-(3-((R)-1-Phenylethoxy)phenyl)-3-(3-chloro-methylbenzo[b]thiophen-2-yl-carbonyl)thiourea;

1-(3-((R)-1-Phenylethoxy)phenyl)-3-(4-trifluoromethoxy-phenyl-carbonyl)thiourea;

1-(3-((R)-1-Phenylethoxy)phenyl)-3-(5-methylisoxazol-3-yl-carbonyl)thiourea;

1-(3-((S)-1-Phenylethoxy)phenyl)-3-(4-trifluoromethylphenyl-carbonyl)-thiourea;

1-(3-((S)-1-Phenylethoxy)phenyl)-3-(4-trifluoromethoxyphenyl-carbonyl)-thiourea;

1-(3-((S)-1-Phenylethoxy)phenyl)-3-(5-methylisoxazol-3-yl-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(4-trifluoromethylphenyl-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(3-chloro-methylbenzo[b]thiophen-2-yl-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(4-trifluoromethoxyphenyl-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(3,5-dimethylisoxazol-4-yl-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(5-methylisoxazol-3-yl-carbonyl)-thiourea;

1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(4-trifluoromethylphenyl-carbonyl)thiourea;

1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(4-trifluoromethoxyphenyl-carbonyl)thiourea;

1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(3,5-dimethylisoxazol-4-yl-carbonyl)thiourea;

1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(5-methylisoxazol-3-yl-carbonyl)thiourea;

1-(2-Phenylbenzo[*d*][1,3]dioxol-6-yl)-3-(benzofuran-2-yl-carbonyl)thiourea;

1-(3-((S)-1-Phenylethoxy)phenyl)-3-(benzofuran-2-yl-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(benzofuran-2-yl-carbonyl)-thiourea;

1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(benzofuran-2-yl-carbonyl)thiourea;

1-(3-((R)-1-Phenylethoxy)phenyl)-3-(benzofuran-2-yl-carbonyl)thiourea;

1-(2,4-dimethylthiazol-5-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(1-methyl-pyrrol-2-yl-carbonyl)-3-(3-(benzyloxy)phenyl)thiourea;

1-(3-(Phenethyloxy)phenyl)-3-((2,7-dimethylpyrazolo[1,5-*a*]pyrimidin-6-yl)-carbonyl)-thiourea;

1-(3-(Phenethyloxy)phenyl)-3-(1-ethyl-3-methyl-1*H*-pyrazol-5-yl-carbonyl)-thiourea;

1-(3-(2,3-Dihydro-1*H*-inden-2-yloxy)phenyl)-3-(2,4-dimethylthiazol-5-yl-carbonyl)thiourea;

1-(3-(2,3-Dihydro-1*H*-inden-2-yloxy)phenyl)-3-((2,7-dimethylpyrazolo[1,5-*a*]pyrimidin-6-yl)-carbonyl)thiourea;

1-(3-(2,3-Dihydro-1*H*-inden-2-yloxy)phenyl)-3-(1-ethyl-3-methyl-1*H*-pyrazol-5-yl-carbonyl)thiourea;

1-(4-Methyl-1,2,3-thiadiazol-5-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(Benzo[*d*]thiazol-2-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;

1-(4-Methyl-1,2,3-thiadiazol-5-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(Benzo[*d*]thiazol-2-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-fluoro-5-pentoxy-phenyl)thiourea;

1-(2-Methyl-pyridin-3-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(1-phenyl-1*H*-pyrazol-5-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(1-phenyl-1*H*-pyrazol-5-yl)-3-((3-phenyloxy)phenyl)-phenyl)thiourea;

1-(2-Phenylbenzo[*d*][1,3]dioxol-6-yl)-3-(1-phenyl-1*H*-pyrazol-5-yl-carbonyl)thiourea;

1-(1-phenyl-1*H*-pyrazol-5-yl)-3-(4-pentoxy-phenyl)-phenyl)thiourea;

1-(1-phenyl-1*H*-pyrazol-5-yl)-3-((3-phenyloxy-phenyl)-phenyl)thiourea;

1-(Methylbenzo[*b*]thiophen-2-yl-carbonyl)-3-((4-pentoxy-phenyl)-phenyl)thiourea;

1-(isoxazol-5-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(isoxazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(isoxazol-5-yl-carbonyl)-3-(4-(pentyl) phenyl)thiourea;

1-(isoxazol-5-yl-carbonyl)-3-(4-(pentoxy) phenyl)thiourea;

1-(isoxazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(isoxazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-((3-(trifluoromethyl)phenyl)furan-2-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-((3-(trifluoromethyl)phenyl)furan-2-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(5-Bromo-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(5-Bromo-benzofuran-2-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(5-Nitro-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(5-Nitro-benzofuran-2-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(5-Bromo-benzofuran-2-yl-carbonyl)-3-(4-pentyl phenyl)thiourea;

1-(5-Bromo-benzofuran-2-yl-carbonyl)-3-(4-pentoxy phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(4-(hex-1-ynyl)phenyl) thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(2-(pyridin-3-yl)ethynyl)phenyl) thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(2-(pyridin-3-yl)ethyl)phenyl) thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(pyridin-2-yl) phenyl) thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(pyridin-2-yl-oxy)-phenyl) thiourea;

1-(3,5-Dimethylisoxazole-4-yl-carbonyl)-3-(4-(hex-1-ynyl)phenyl) thiourea;

1-(3,5-Dimethylisoxazole-4-yl-carbonyl)-3-(3-(2-(pyridin-3-yl)ethynyl)phenyl) thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-((3-(trifluoromethyl)benzyloxy)-phenyl) thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(4-((1-methylpiperidin-4-yl)methoxy)-3-fluorophenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(trifluoromethyl)-4-(piperidin-1-yl)phenyl)thiourea;

1-(1,3-dimethyl-1*H*-thieno[2,3-*c*]pyrazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(5-(2-methylthiazol-4-yl)isoxazol-3-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(5-(2-methylthiazol-4-yl)isoxazol-3-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(5-(2-methylthiazol-4-yl)isoxazol-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(5-(2-methylthiazol-4-yl)isoxazol-3-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(1,5-Dimethyl-1*H*-pyrazol-3-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(1,5-Dimethyl-1*H*-pyrazol-3-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(1,5-Dimethyl-1*H*-pyrazol-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(1,5-Dimethyl-1*H*-pyrazol-3-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(1-Methyl-3-*tert*-butyl-1*H*-pyrazol-3-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(1-Methyl-3-*tert*-butyl-1*H*-pyrazol-3-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(1-Methyl-3-*tert*-butyl-1*H*-pyrazol-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(2-Trifluoromethyl-5-methyl-furan-2-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(2-Trifluoromethyl-5-methyl-furan-2-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(2-Trifluoromethyl-5-methyl-furan-2-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(Benzo[*c*][1,2,5]oxadiazol-5-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(Benzo[*c*][1,2,5]oxadiazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;

1-(Benzo[*c*][1,2,5]oxadiazol-5-yl-carbonyl)-3-(4-(pentyl) phenyl)thiourea;

1-(Benzo[*c*][1,2,5]oxadiazol-5-yl-carbonyl)-3-(4-(pentoxy) phenyl)thiourea;

1-(2,7-Dimethylpyrazolo[1,5-*a*]pyrimidin-6-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(2,7-Dimethylpyrazolo[1,5-*a*]pyrimidin-6-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(2,7-Dimethylpyrazolo[1,5-*a*]pyrimidin-6-yl-carbonyl)-3-(4-(pentyl)-phenyl)thiourea;

1-(3-Methylisoxazol-4-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(3-Methylisoxazol-4-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(3-Methylisoxazol-4-yl-carbonyl)-3-(4-(pentyl)-phenyl)thiourea;

1-(3-Methylisoxazol-4-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(5-Methyl-2-phenyl-2*H*-1,2,3-triazol-4-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(5-Methyl-2-phenyl-2*H*-1,2,3-triazol-4-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(5-Chloro-3-methylbenzo[*b*]thiophen-2-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(5-Chloro-3-methylbenzo[*b*]thiophen-2-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(1,3-Dimethyl-1*H*-pyrazol-5-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(1,3-Dimethyl-1*H*-pyrazol-5-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(1,3-Dimethyl-1*H*-pyrazol-5-yl-carbonyl)-3-(4-(pentyl)-phenyl)thiourea;

1-(1,3-Dimethyl-1*H*-pyrazol-5-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(2-(Pyridin-3-yl)thiazol-4-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(2-(Pyridin-3-yl)thiazol-4-yl-carbonyl)-3-(3-(phenyloxy)-phenyl)thiourea;

1-(2-(Pyridin-3-yl)thiazol-4-yl-carbonyl)-3-(4-(pentyl)-phenyl)thiourea;

1-(2-(Pyridin-3-yl)thiazol-4-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(4-Methoxy-benzofuran-2-yl-carbonyl)-3-(4-(benzyloxy)-phenyl)thiourea;

1-(4-Methoxy-benzofuran-2-yl-carbonyl)-3-(3-fluoro-4-pentoxy-phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3,5-dibromo-4-(pent-4-enyloxy)phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-((pyridin-3-yl)methyl) phenyl)thiourea;

1-(5-Iodo-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)phenyl)thiourea;

1-(5-Phenyl-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)phenyl)thiourea;

1-(5-(2-Pyridyl)benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)phenyl)thiourea;

1-(3-Propoxy-pyridin-2-yl-carbonyl)-3-(4-(pentyl)phenyl)thiourea;

1-(2,5-Dichlorothiophen-3-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(3-Methyl-5-(methylthio)-4-vinylthiophen-2-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(3-Methyl-5-(methylthio)-4-vinylthiophen-2-yl-carbonyl)-3-(4-(benzyloxy)-phenyl)thiourea;

1-(5-(Methylthio)-thiophen-2-yl-carbonyl)-3-(4-(benzyloxy)-phenyl)thiourea;

1-(5,7-Dimethylpyrazolo[1,5-*a*]pyrimidin-2-yl-carbonyl)-3-(4-(benzyloxy)-phenyl)thiourea;

1-(2,5-Dichlorothiophen-3-yl-carbonyl)-3-(4-(phenoxy)-phenyl)thiourea;

1-(3-Methyl-5-(methylthio)-4-vinylthiophen-2-yl-carbonyl)-3-(4-(phenoxy)-phenyl)thiourea;

1-(5-(Methylthio)-thiophen-2-yl-carbonyl)-3-(4-(phenoxy)-phenyl)thiourea;

1-(7-Fluoro-benzofuran-2-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(7-Fluoro-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(2-(1,3-Dioxoisoindolin-2-yl)acetyl)-3-(3-phenoxyphenyl)thiourea;

1-(2-(1,3-Dioxoisoindolin-2-yl)acetyl)-3-(3-benzyloxyphenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)methyl-phenyl)thiourea ;

1-(Benzofuran-2-yl-carbonyl)-3-(3-(phenylamino)methyl-phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(4-(*N*-benzyl-N-methylamino)-3-fluorophenyl) thiourea;

Phenyl 3-(3-((benzofuran-2-yl)-carbonyl)thioureido)benzoate;

1-(4-Cyanophenyl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(4-Cyanophenyl-carbonyl)-3-(4-(pentyloxy)-phenyl)thiourea;

1-(4-Cyanophenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(Quinoxalin-2-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(Quinoxalin-2-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(Quinoxalin-2-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(4-Trifluoromethoxyphenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(4-Trifluoromethoxyphenyl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(4-Trifluoromethylphenyl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(4-Trifluoromethylphenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(5-Cyano-benzofuran-2-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(5-Cyano-benzofuran-2-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(4-Methyl-1,2,3-thiadiazol-5-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(4-Methyl-1,2,3-thiadiazol-5-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

Butyl 4-(3-(4-Methyl-1,2,3-thiadiazol-5-yl-) carbonyl)thioureido) benzoate;

1-(Pyrazin-2-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(Pyrazin-2-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(Pyrazin-2-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(3-Fluoro-4-(pentyloxy)phenyl)-3-(2-(1,3-dioxoisoindolin-2-yl)acetyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(3-trifluoromethyl-4-pentoxy-phenyl)thiourea;

1-(1-Benzyl-1*H*-tetrazol-5-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(4-(N-methyl-N-pentylamino)-3-fluorophenyl)thiourea;

1-(1-Benzyl-1*H*-tetrazol-5-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(6-Trifluoromethyl-pyrid-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(6-Trifluoromethyl-pyrid-3-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(3-Trifluoromethyl-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(3-Trifluoromethyl-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(3-Trifluoromethyl-phenyl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(3-Trifluoromethoxy-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(3-Trifluoromethoxy-phenyl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(2-Chloro-5-trifluoromethoxy-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(3-Difluoromethyl-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(3-Difluoromethyl-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(5-(Trifluoromethyl)-2-phenyloxazol-4-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(5-(Trifluoromethyl)-2-phenyloxazol-4-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;

1-(5-(2-Chloro-5-trifluoromethylphenyl)-furan-2-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;

1-(3-Trifluoromethyl-4-methoxy-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(3-Trifluoromethyl-4-chloro-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(3-Trifluoromethyl-4-chloro-phenyl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(3-Trifluoromethyl-4-methyl-phenyl-carbonyl)-3-(3-benzyloxy-phenyl)thiourea;

1-(3-Trifluoromethyl-4-methyl-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(3-Trifluoromethyl-4-methyl-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-((5-Acetamidobenzofuran-2-yl)carbonyl)-3-(3-phenoxyphenyl)thiourea;

1-Acetyl-3-(3-phenoxyphenyl)thiourea;

1-Acetyl-3-(4-(pentyloxy)phenyl)thiourea;

1-Acetyl-3-(4-pentylphenyl)thiourea;

1-(Dimethylamino-acetyl)-3-(3-phenoxyphenyl)thiourea;

1-(Dimethylamino-acetyl)-3-(3-benzyloxyphenyl)thiourea;

1-(3,5-Dimethylisoxazole-4-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-((2,3,4,5,6-penta-fluorophenoxy)-phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(9-methyl-9*H*-fluoren-7-yl)thiourea;

Pentyl 2-phenyl- 4-(3-(benzofuran-2-yl)thioureido)benzoate;

1-(3-Pyrid-3-yl-carbonyl)-3-(3-benzyloxy-phenyl)thiourea;

1-(3-Pyrid-3-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;

1-(3-Pyrid-3-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;

1-(3-Pyrid-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(4-Phenylbutanoyl)-3-(3-phenoxyphenyl)thiourea;

1-(4-Phenylbutanoyl)-3-(3-benzyloxyphenyl)thiourea;

1-(2-Morpholinoacetyl)-3-(3-phenoxyphenyl)thiourea;

1-(2-Morpholinoacetyl)-3-(4-(pentyloxy)phenyl)thiourea;

1-(2-Morpholinoacetyl)-3-(4-(pentyl)phenyl)thiourea;

1-(4-(Pentyloxy)phenyl)-3-(2-(piperidin-1-yl)acetyl)thiourea;

1-(*N*-Methyl-*N*-phenylamino-acetyl)-3-(3-benzyloxyphenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(6-pentoxy-pyrid-3-yl)thiourea;

1-(3-Pyrid-3-yl-carbonyl)-3-(3-benzyloxy-phenyl)thiourea hydrochloride;

1-(3-Pyrid-3-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea hydrochloride;

1-(3-Pyrid-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea hydrochloride;

1-(Benzofuran-2-yl-carbonyl)-3-(trifluoromethylthio-phenyl)thiourea;

1-(3-(Piperidin-1-yl)propanoyl)-3-(4-pentylphenyl)thiourea;

1-(3-(Piperidin-1-yl)propanoyl)-3-(4-(pentyloxy)phenyl)thiourea;

1-(3-(Piperidin-1-yl)propanoyl)-3-(3-phenoxyphenyl)thiourea;

1-(3-Morpholinopropanoyl)-3-(4-(pentyloxy)phenyl)thiourea;

1-(1-Methylpiperidin-3-yl-carbonyl)-3-(4-(pentyloxy)phenyl)thiourea;

1-(1-Methylpiperidin-3-yl-carbonyl)-3-(4-(pentyloxy)phenyl)thiourea;

1-(2-(2-methylpiperidin-1-yl)acetyl)-3-(4-(pentyloxy)phenyl)thiourea;

1-(2-Oxo-4-phenyl-pyrrolidin-1-ylcarbonyl)-3-(3-benzyloxy-phenyl)thiourea; and

1-(5-Trifluoromethoxy-benzofuran-2-yl-carbonyl)-3-(3-benzyloxy-phenyl)thiourea.

82. (Currently Amended) A pharmaceutical composition comprising a compound or salt according to ~~any one of Claims 1 to 81~~ Claim 1 together with a pharmaceutically acceptable carrier, diluent, or excipient.

Claims 83-86. (Canceled)

87. (Original) A method for treating Hepatitis C infection comprising administering to a patient in need of such treatment a therapeutically effective amount of a compound or salt according to Claim 1.

Claims 88-90. (Canceled)